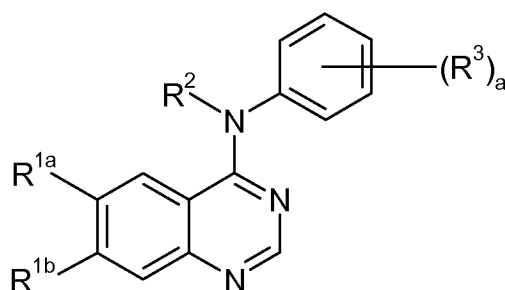


CLAIM AMENDMENTS:

This listing of claims will replace all prior versions and listing of claims in the application.

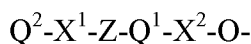
Listing of the Claims:

Claim 1 (**currently amended**): A quinazoline derivative of the Formula I:



wherein:

one of R^{1a} or R^{1b} is a group of sub-formula (i)



(i)

where X² and X¹ are independently selected from a direct bond or a group -[CR⁴R⁵]_m, wherein m is an integer from 1 to 6,

Z is C(O), SO₂, -C(O)NR¹⁰-, -N(R¹⁰)C(O)-, -C(O)O- or -OC(O)- where R¹⁰ is hydrogen or (1-6C)alkyl,

and each of R⁴ and R⁵ is independently selected from hydrogen, hydroxy, (1-4C)alkyl, halo(1-4C)alkyl, hydroxy (1-4C)alkyl, (1-4C)alkoxy(1-4C)alkyl, or R⁴ and R⁵ together with the carbon atom(s) to which they are attached form a (3-7)cycloalkyl ring, provided that when a group R⁴ or R⁵ is hydroxy, m is at least 2 and the carbon atom to which the hydroxy group is attached is not also attached to another oxygen or a nitrogen atom;

Q¹ is a piperidinyl ring, which is optionally substituted by one or two substituents selected from halogeno, trifluoromethyl, trifluoromethoxy, cyano, nitro, hydroxy, amino, carboxy, carbamoyl,

acryloyl, (1-6C)alkyl, (2-8C)alkenyl, (2-8C)alkynyl, (1-6C)alkoxy, (2-6C)alkenyloxy, (2-6C)alkynyloxy, (1-6C)alkylthio, (2-6C)alkenylthio, (2-6C)alkynylthio, (1-6C)alkylsulfinyl, (2-6C)alkenylsulfinyl, (2-6C)alkynylsulfinyl, (1-6C)alkylsulfonyl, (2-6C)alkenylsulfonyl, (2-6C)alkynylsulfonyl, (1-6C)alkylamino, di-[(1-6C)alkyl]amino, (1-6C)alkoxycarbonyl, N-(1-6C)alkylcarbamoyl, N,N-di-[(1-6C)alkyl]carbamoyl, (2-6C)alkanoyl, (2-6C)alkanoyloxy, (2-6C)alkanoylamino, N-(1-6C)alkyl-(2-6C)alkanoylamino, sulfamoyl, N-(1-6C)alkylsulfamoyl, N,N-di-[(1-6C)alkyl]sulfamoyl, (1-6C)alkanesulfonylamino, N-(1-6C)alkyl-(1-6C)alkanesulfonylamino, carbamoyl(1-6C)alkyl, N-(1-6C)alkylcarbamoyl(1-6C)alkyl, N,N-di-[(1-6C)alkyl]carbamoyl(1-6C)alkyl, sulfamoyl(1-6C)alkyl, N-(1-6C)alkylsulfamoyl(1-6C)alkyl, N,N-di-[(1-6C)alkyl]sulfamoyl(1-6C)alkyl, (2-6C)alkanoyl(1-6C)alkyl, (2-6C)alkanoyloxy(1-6C)alkyl, (2-6C)alkanoylamino(1-6C)alkyl, N-(1-6C)alkyl-(2-6C)alkanoylamino(1-6C)alkyl and (1-6C)alkoxycarbonyl(1-6C)alkyl;

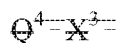
Q² is an isoxazolyl ring optionally substituted by one or two groups, which may be the same or different, selected from halogeno, hydroxy, nitro, amino, cyano, carbamoyl, (1-4C)alkyl, (1-4C)alkoxy, (2-4C)alkanoyl and (1-4C)alkylsulfonyl, [(1-4C)alkyl]amino, di[(1-4C)alkyl]amino, *N*-[(1-4C)alkyl]carbamoyl, and *N,N*-di[(1-4C)alkyl]carbamoyl;

and wherein any (2-4C)alkanoyl group in a substituent on Q² optionally bears one or two substituents, which may be the same or different, selected from hydroxy and (1-3C)alkyl,

and wherein any (1-4C)alkyl group in a substituent on Q² optionally bears one or two substituents, which may be the same or different, selected from hydroxy, (1-4C)alkoxy and halogeno; ~~one of more substituents selected from halogeno, trifluoromethyl, trifluoromethoxy, cyano, nitro, hydroxy, amino, carboxy, carbamoyl, acryloyl, (1-6C)alkyl, (2-8C)alkenyl, (2-8C)alkynyl, (1-6C)alkoxy, (2-6C)alkenyloxy, (2-6C)alkynyloxy, (1-6C)alkylthio, (2-6C)alkenylthio, (2-6C)alkynylthio, (1-6C)alkylsulfinyl, (2-6C)alkenylsulfinyl, (2-6C)alkynylsulfinyl, (1-6C)alkylsulfonyl, (2-6C)alkenylsulfonyl, (2-6C)alkynylsulfonyl, (1-6C)alkylamino, di-[(1-6C)alkyl]amino, (1-6C)alkoxycarbonyl, N-(1-6C)alkylcarbamoyl, N,N-di-[(1-6C)alkyl]carbamoyl, (2-6C)alkanoyl, (2-6C)alkanoyloxy, (2-6C)alkanoylamino, N-(1-6C)alkyl-(2-6C)alkanoylamino, sulfamoyl, N-(1-6C)alkylsulfamoyl,~~

~~N,N-di-[(1-6C)alkyl]sulfamoyl, (1-6C)alkanesulfonylamino,
N-(1-6C)alkyl-(1-6C)alkanesulfonylamino, carbamoyl(1-6C)alkyl,
N-(1-6C)alkylcarbamoyl(1-6C)alkyl, N,N-di-[(1-6C)alkyl]carbamoyl(1-6C)alkyl,
sulfamoyl(1-6C)alkyl, N-(1-6C)alkylsulfamoyl(1-6C)alkyl,
N,N-di-[(1-6C)alkyl]sulfamoyl(1-6C)alkyl, (2-6C)alkanoyl(1-6C)alkyl,
(2-6C)alkanoyloxy(1-6C)alkyl, (2-6C)alkanoylamino(1-6C)alkyl,
N-(1-6C)alkyl-(2-6C)alkanoylamino(1-6C)alkyl and (1-6C)alkoxy-carbonyl(1-6C)alkyl,
and wherein any (1-6C)alkyl, (2-8C)alkenyl, (2-8C)alkynyl and (2-6C)alkanoyl
substituent on Q¹ or Q² optionally bears one or more substituents which may be the same or
different selected from halogeno, hydroxy and (1-6C)alkyl and/or optionally a substituent
selected from cyano, nitro, carboxy, (2-8C)alkenyl, (2-8C)alkynyl, (1-6C)alkoxy,
hydroxy(1-6C)alkoxy, (1-4C)alkoxy(1-6C)alkoxy, (2-6C)alkanoyl, (2-6C)alkanoyloxy and
NR^aR^b, wherein R^a is hydrogen or (1-4C)alkyl and R^b is hydrogen or (1-4C)alkyl, and wherein
any (1-4C)alkyl in R^a or R^b optionally bears one or more substituents which may be the same or
different selected from halogeno and hydroxy and/or optionally a substituent selected from
cyano, nitro, (2-4C)alkenyl, (2-4C)alkynyl, (1-4C)alkoxy, hydroxy(1-4C)alkoxy and
(1-2C)alkoxy(1-4C)alkoxy,
or R^a and R^b together with the nitrogen atom to which they are attached form a 4, 5 or 6
membered ring, which optionally bears 1 or 2 substituents, which may be the same or different,
on an available ring carbon atom selected from halogeno, hydroxy, (1-4C)alkyl and
(1-3C)alkylenedioxy, and may optionally bear on any available ring nitrogen a substituent
(provided the ring is not thereby quaternised) selected from (1-4C)alkyl, (2-4C)alkanoyl and
(1-4C)alkylsulfonyl,
and wherein any (1-4C)alkyl or (2-4C)alkanoyl group present as a substituent on the ring
formed by R^a and R^b together with the nitrogen atom to which they are attached, optionally bears
one or more substituents which may be the same or different selected from halogeno and
hydroxy and/or optionally a substituent selected from (1-4C)alkyl and (1-4C)alkoxy;
and wherein Q¹ optionally bears 1 or 2 oxo (=O) or thioxo (=S) substituents;~~

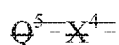
and the other of R^{1a} or R^{1b} is a group R^1 which is hydrogen, (1-6C)alkoxy and (1-4C)alkoxy(1-6C)alkoxy, and wherein any (1-6C)alkoxy group within R^1 optionally bears 1, 2 or 3 substituents, which may be the same or different, selected from hydroxy, fluoro and chloro~~selected from hydrogen, hydroxy, (1-6C)alkoxy, (2-6C)alkenyloxy, (2-6C)alkynyloxy, or a group of the formula:~~



wherein X^3 is a direct bond or is selected from O or S, and Q^4 is (3-7C)cycloalkyl, (3-7C)cycloalkyl-(1-6C)alkyl, (3-7C)cycloalkenyl, (3-7C)cycloalkenyl-(1-6C)alkyl, heterocycetyl or heterocycetyl-(1-6C)alkyl,

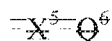
~~and wherein adjacent carbon atoms in any (2-6C)alkylene chain within a R^+ substituent are optionally separated by the insertion into the chain of a group selected from O, S, SO, SO₂, N(R^4), CO, CH(OR⁴), CON(R^4), N(R^4)CO, SO₂N(R^4), N(R^4)SO₂, CH=CH and C≡C wherein R^4 is hydrogen or (1-6C)alkyl,~~

~~and wherein any CH₂=CH or HC≡C group within a R^+ substituent optionally bears at the terminal CH₂= or HC≡ position a substituent selected from halogeno, carboxy, carbamoyl, (1-6C)alkoxycarbonyl, N-(1-6C)alkylcarbamoyl, N,N-di-[(1-6C)alkyl]carbamoyl, amino-(1-6C)alkyl, (1-6C)alkylamino-(1-6C)alkyl and di-[(1-6C)alkyl]amino-(1-6C)alkyl or from a group of the formula:~~



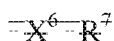
wherein X^4 is a direct bond or is selected from CO and N(R^5)CO, wherein R^5 is hydrogen or (1-6C)alkyl, and Q^5 is heterocycetyl or heterocycetyl-(1-6C)alkyl,

~~and wherein any alkyl or alkylene group within a R^+ substituent optionally bears one or more halogeno, (1-6C)alkyl, hydroxy, cyano, amino, carboxy, carbamoyl, sulfamoyl, (1-6C)alkoxy, (1-6C)alkylthio, (1-6C)alkylsulfinyl, (1-6C)alkylsulfonyl, (1-6C)alkylamino, di-[(1-6C)alkyl]amino, (1-6C)alkoxycarbonyl, N-(1-6C)alkylcarbamoyl, N,N-di-[(1-6C)alkyl]carbamoyl, (2-6C)alkanoyl, (2-6C)alkanoyloxy, (2-6C)alkanoylamino, N-(1-6C)alkyl-(2-6C)alkanoylamino, N-(1-6C)alkylsulfamoyl, N,N-di-[(1-6C)alkyl]sulfamoyl, (1-6C)alkanesulfonylamino and N-(1-6C)alkyl-(1-6C)alkanesulfonylamino, or from a group of the formula:~~



wherein X^5 is a direct bond or is selected from O, S, SO, SO₂, N(R⁶), CO, CH(OR⁶), CON(R⁶), N(R⁶)CO, SO₂N(R⁶), N(R⁶)SO₂, C(R⁶)₂O, C(R⁶)₂S and C(R⁶)₂N(R⁶), wherein R⁶ is hydrogen or (1-6C)alkyl, and Q⁶ is (3-7C)cycloalkyl, (3-7C)cycloalkyl-(1-6C)alkyl, (3-7C)cycloalkenyl, (3-7C)cycloalkenyl-(1-6C)alkyl, heterocycetyl or heterocycetyl-(1-6C)alkyl;

and wherein any heterocycetyl group within a substituent on R¹ optionally bears 1, 2 or 3 substituents, which may be the same or different, selected from halogeno, trifluoromethyl, cyano, nitro, hydroxy, amino, carboxy, carbamoyl, formyl, mercapto, (1-6C)alkyl, (2-8C)alkenyl, (2-8C)alkynyl, (1-6C)alkoxy, (2-6C)alkenyloxy, (2-6C)alkynyloxy, (1-6C)alkylthio, (1-6C)alkylsulfinyl, (1-6C)alkylsulfonyl, (1-6C)alkylamino, di-[(1-6C)alkyl]amino, (1-6C)alkoxycarbonyl, N-(1-6C)alkylcarbamoyl, N,N-di-[(1-6C)alkyl]carbamoyl, (2-6C)alkanoyl, (2-6C)alkanoyloxy, (2-6C)alkanoylamino, N-(1-6C)alkyl-(2-6C)alkanoylamino, N-(1-6C)alkylsulfamoyl, N,N-di-[(1-6C)alkyl]sulfamoyl, (1-6C)alkanesulfonylamino, and N-(1-6C)alkyl-(1-6C)alkanesulfonylamino, or from a group of the formula:



wherein X^6 is a direct bond or is selected from O, N(R⁸) and C(O), wherein R⁸ is hydrogen or (1-6C)alkyl, and R⁷ is halogeno (1-6C)alkyl, hydroxy (1-6C)alkyl, carboxy (1-6C)alkyl, (1-6C)alkoxy (1-6C)alkyl, cyano (1-6C)alkyl, amino (1-6C)alkyl, (1-6C)alkylamino (1-6C)alkyl, di-[(1-6C)alkyl]amino (1-6C)alkyl, (2-6C)alkanoylamino (1-6C)alkyl, (1-6C)alkoxycarbonylamino (1-6C)alkyl, carbamoyl (1-6C)alkyl, N-(1-6C)alkylcarbamoyl (1-6C)alkyl, N,N-di-[(1-6C)alkyl]carbamoyl (1-6C)alkyl, (2-6C)alkanoyl (1-6C)alkyl or (1-6C)alkoxycarbonyl (1-6C)alkyl;

and wherein any heterocycetyl group within a substituent on R¹ optionally bears 1 or 2 exo or thio substituents;

R² is selected from hydrogen and (1-6C)alkyl;

each R³, which may be the same or different, is selected from halogeno, cyano, nitro,

hydroxy, amino, carboxy, carbamoyl, sulfamoyl, trifluoromethyl, (1-6C)alkyl, (2-8C)alkenyl, (2-8C)alkynyl, (1-6C)alkoxy, (2-6C)alkenyloxy, (2-6C)alkynyloxy, (1-6C)alkylthio, (1-6C)alkylsulfinyl, (1-6C)alkylsulfonyl, (1-6C)alkylamino, di-[(1-6C)alkyl]amino, (1-6C)alkoxycarbonyl, N-(1-6C)alkylcarbamoyl, N,N-di-[(1-6C)alkyl]carbamoyl, N-(1-6C)alkylsulfamoyl, and N,N-di-[(1-6C)alkyl]sulfamoyl

a is 1, 2 or ~~3, 3, 4 or 5~~;

or a pharmaceutically acceptable salt thereof;

~~subject to the proviso that the compound of formula I is not~~

~~N-(3,4-dichlorophenyl)-7-[(4-[(3,5-dimethylisoxazol-4-yl)carbonyl]morpholin-2-yl)methoxy]-6-(methoxy)quinazolin-4-amine.~~

Claim 2 (**previously presented**): The quinazoline derivative according to claim 1 wherein X² is a direct bond.

Claims 3-6 (**cancelled**).

Claim 7 (**currently amended**): The quinazoline derivative according to claim ~~1-6~~ wherein R¹ is selected from methoxy, ethoxy, isopropoxy, cyclopropylmethoxy, 2-hydroxyethoxy, 2-fluoroethoxy, 2-methoxyethoxy, 2,2-difluoroethoxy, 2,2,2-trifluoroethoxy or 3-hydroxy-3-methylbutoxy.

Claim 8 (**currently amended**): The quinazoline derivative according to claim ~~7-5~~ wherein R¹ is methoxy.

Claim 9 (**previously presented**): The quinazoline derivative according to claim 1 wherein X¹ is suitably a direct bond or a (1-6C)alkylene group.

Claim 10 (**previously presented**): The quinazoline derivative according to claim 9 wherein X¹ is a direct bond or methylene or ethylene group.

Claim 11 (**previously presented**): The quinazoline derivative according to claim 1 wherein Z is selected from -C(O)-, -NR¹⁰-C(O)- (wherein R¹⁰ is H or (1-6C)alkyl), and -O-C(O)-.

Claim 12 (**previously presented**): The quinazoline derivative according to claim 11, wherein Z is -C(O)-.

Claim 13 (**previously presented**): The quinazoline derivative according to claim 11, wherein Z is selected from -NH-C(O)- and -O-C(O)-.

Claims 14-15 (**cancelled**).

Claim 16 (**currently amended**): The quinazoline derivative according to claim ~~1-11~~, wherein the group Q²-X¹-Z- is linked to the piperidinyl nitrogen of Q¹.

Claims 17-24 (**cancelled**).

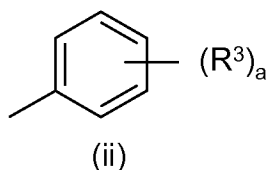
Claim 25 (**currently amended**): The quinazoline derivative according to claim ~~1-23~~ wherein Q² is unsubstituted or substituted by a (1-4C)alkyl group, a (1-4C)alkoxy group, halogeno, amino, nitro, cyano, carbamoyl, di-[(1-4C)alkyl]amino, and *N,N*-di[(1-4C)alkyl]carbamoyl.

Claim 26 (**previously presented**): The quinazoline derivative according to claim 1 wherein R² is hydrogen.

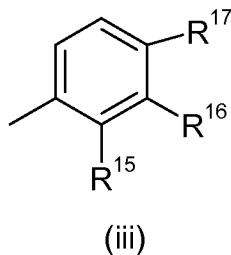
Claim 27 (**cancelled**).

Claim 28 (**previously presented**): The quinazoline derivative according to claim 1, wherein an R³ is in the para position on the anilino ring, and this is selected from halogeno, cyano, nitro, hydroxy, amino, trifluoromethyl, (1-6C)alkyl, (2-8C)alkenyl, (2-8C)alkynyl, (1-6C)alkoxy, (2-6C)alkenyloxy, (2-6C)alkynyloxy, (1-6C)alkylthio, (1-6C)alkylamino and di-[(1-6C)alkyl]amino.

Claim 29 (**previously presented**): The quinazoline derivative according to claim 1 wherein the group of sub-formula (ii)



in formula (I) is a group of sub-formula (iii)



where one of R¹⁵ or R¹⁷ is hydrogen and the other is halogeno, and R¹⁶ is halogeno.

Claim 30 (**previously presented**): The quinazoline derivative according to claim 29 wherein the group of sub-formula (iii) is 3-chloro-2-fluorophenyl, or 3-chloro-4-fluorophenyl.

Claim 31 (**currently amended**): The compound according to claim 1 selected from one of the following:

- (1) *N*-(3-chloro-2-fluorophenyl)-6-{[1-(isoxazol-5-ylcarbonyl)piperidin-4-yl]oxy}-7-methoxyquinazolin-4-amine;
- (2) *N*-(3-chloro-2-fluorophenyl)-7-methoxy-6-({ 1-[(3-methylisoxazol-5-yl)acetyl]piperidin-4-yl } oxy)quinazolin-4-amine;

- (3) *N*-(3-chloro-2-fluorophenyl)-7-methoxy-6-({1-[(3-methylisoxazol-5-yl)carbonyl]piperidin-4-yl}oxy)quinazolin-4-amine;
- (4) *N*-(3-chloro-2-fluorophenyl)-7-methoxy-6-({1-[(5-methylisoxazol-3-yl)carbonyl]piperidin-4-yl}oxy)quinazolin-4-amine;
- (5) *N*-(3-chloro-2-fluorophenyl)-7-methoxy-6-({1-[(5-methylisoxazol-4-yl)carbonyl]piperidin-4-yl}oxy)quinazolin-4-amine;
- (6) *N*-(3-chloro-2-fluorophenyl)-7-methoxy-6-({1-[(3-methylisoxazol-4-yl)carbonyl]piperidin-4-yl}oxy)quinazolin-4-amine;
- (7) *N*-(3-chloro-2-fluorophenyl)-6-({1-[(3,5-dimethylisoxazol-4-yl)carbonyl]piperidin-4-yl}oxy)-7-methoxyquinazolin-4-amine;
- (8) *N*-(3-chloro-2-fluorophenyl)-7-{{1-[(isoxazol-5-ylcarbonyl)piperidin-4-yl]oxy}-6-methoxyquinazolin-4-amine;
- (9) *N*-(3-chloro-2-fluorophenyl)-6-methoxy-7-({1-[(3-methylisoxazol-5-yl)acetyl]piperidin-4-yl}oxy)quinazolin-4-amine;
- (10) *N*-(3-chloro-2-fluorophenyl)-7-methoxy-6-({(3*R*)-1-[(3-methylisoxazol-5-yl)acetyl]piperidin-3-yl}oxy)quinazolin-4-amine; and
- ~~(11) *N*-(3-chloro-2-fluorophenyl)-7-methoxy-6-{{(3*R*)-1-(4-{*N,N*-dimethylcarbamoyl}-1*H*-pyrazol-1-yl)acetyl}piperidin-3-yl}oxy}quinazolin-4-amine; and~~
- ~~(12)~~ (11) 4-({4-[(3-Chloro-2-fluorophenyl)amino]-7-methoxyquinazolin-6-yl}oxy)-*N*-(3,5-dimethylisoxazol-4-yl)piperidine-1-carboxamide.

Claims 32-33 (**cancelled**).

Claim 34 (**previously presented**): A pharmaceutical composition which comprises a quinazoline derivative of the Formula I, or a pharmaceutically-acceptable salt thereof, as defined in claim 1 in association with a pharmaceutically-acceptable diluent or carrier.

Claims 35-37 (**cancelled**).